

# An asymptotic approximation for the permanent of a doubly stochastic matrix

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## Abstract

A determinantal approximation is obtained for the permanent of a doubly stochastic matrix. For moderate-deviation matrix sequences, the asymptotic relative error is of order  $O(n^{-1})$ .

**keywords:** Doubly stochastic Dirichlet distribution; Maximum-likelihood projection; Sinkhorn projection

## 1 Introduction

A non-negative matrix of order  $n$  with unit row and column sums is called doubly stochastic. To each strictly positive square matrix  $Y$  there corresponds a unique re-scaled matrix  $A$  such that  $Y_{ij} = nA_{ij} \exp(\beta_i + \gamma_j)$  where  $A$  is strictly positive with unit row and column sums. The doubly stochastic projection  $Y \mapsto A$  can be computed by iterative proportional scaling (Deming and Stephan, 1940), although, in this context, it is usually called the Sinkhorn algorithm (Sinkhorn, 1964; Linial, Samorodnitsky and Wigderson, 1998). A non-negative matrix  $Y$  containing some zero components is said to be scalable if there exist real numbers  $\beta_i, \gamma_j$  such that  $Y_{ij} = nA_{ij} \exp(\alpha_i + \gamma_j)$  where  $A$  is doubly stochastic. In that case, the pattern of zeros in  $Y$  is the same as the pattern of zeros in  $A$ . A non-negative matrix is scalable if and only if  $\text{per}(Y) > 0$ .

The permanent of a square matrix  $A$

$$\text{per}(A) = \sum_{\sigma} \prod_{i=1}^n A_{i,\sigma(i)}$$

is the sum over permutations  $\sigma: [n] \rightarrow [n]$  of products,  $A_{1,\sigma(1)} \cdots A_{n,\sigma(n)}$ , one component taken from each row and each column. If  $A_{ij} = \alpha_i \beta_j B_{ij}$ , then the permanents are related by

$$\text{per}(A) = \text{per}(B) \prod_{i=1}^n (\alpha_i \beta_j).$$

Thus, the problem of approximating the permanent of a scalable non-negative matrix is reduced to the approximation of the corresponding doubly stochastic matrix, which is the subject of this paper.

It is known that the permanent of a generic matrix is not computable in polynomial time as a function of  $n$  (Valiant, 1979). Since exact computation is not feasible for large matrices, most authors have sought to approximate the permanent using Monte-Carlo methods. For example, Jerrum and Sinclair (1989) and Jerrum, Sinclair and Vigoda (2004) use a Markov chain technique, while Kou and McCullagh (2009) use an importance-sampling scheme for  $\alpha$ -permanent approximation. This paper develops an entirely different sort of computable approximation, deterministic but asymptotic.

The determinantal formula (1) is an asymptotic approximation for the permanent of a doubly stochastic matrix  $B$ , the conditions for which are ordinarily satisfied if the original matrix  $A$  has independent and identically distributed components. The approximation is not universally valid for all matrix sequences, but it is valid under moderate-deviation conditions, which can be checked. For example,  $A$  may be symmetric.

## 2 Sinkhorn projection

The scaling algorithm has an interpretation connected with maximum likelihood for generalized linear models as follows. Let the components of  $Y$  be independent exponential variables with means  $\mu_{ij}$  such that  $\log \mu_{ij} = \beta_i + \gamma_j$  lies in the additive subspace *row+col*. The maximum-likelihood projection  $Y \mapsto \hat{\mu}$  is such that

$$\sum_i (Y_{ij}/\hat{\mu}_{ij} - 1) = \sum_j (Y_{ij}/\hat{\mu}_{ij} - 1) = 0$$

(McCullagh and Nelder, 1989). The residual matrix  $Y/\hat{\mu}$  is strictly positive with row and column sums equal to  $n$ , from which it follows that  $A = n^{-1}Y/\hat{\mu}$  is doubly stochastic. The residual deviance

$$\text{Dev}(A) = -2 \sum_{ij} \log(nA_{ij})$$

is a measure of total variability for strictly positive doubly stochastic matrices, taking the value zero only for the uniform matrix  $A_{ij} = 1/n$ . The exponential assumption can be replaced by any gamma distribution provided that the gamma index  $\nu$  is held constant. The re-scaled matrix  $A$  thus generated is said to have the doubly stochastic Dirichlet distribution  $\text{DSD}_n(\nu)$  with parameter  $\nu$ . The smaller the value of  $\nu$ , the more extreme the components of  $A$ , and the greater the deviance. The distribution  $\text{DSD}_n(1)$  is not uniform with respect to Lebesgue measure in the sense of Chatterjee, Diaconis and Sly (2010), but it is presumably close for large  $n$ . The maximum-likelihood estimate  $\hat{\nu}(A)$  is such that

$$2\log(\hat{\nu}) - 2\psi(\hat{\nu}) = \text{Dev}(A)/n^2,$$

where  $\psi$  is the derivative of the log gamma function. It is apparent from simulations that if  $A \sim \text{DSD}_n(\nu)$ ,

$$2\hat{\nu}(A) \times \log(\text{per}(nA)/n!) = 1 + O_p(n^{-1}),$$

with variability decreasing in  $n$ . This limiting product appears to be invariably less than one for non-Dirichlet matrices.

The scaling  $Y \mapsto A$  can be accomplished either by iterative weighted least squares, as is usually done for generalized linear models, or by the Sinkhorn iterative proportional scaling algorithm. For present purposes, the latter algorithm is preferred because it is efficient and simple to implement. Linial, Samorodnitsky and Wigderson (1998) provide a modification that guarantees convergence in polynomial time. For moderate-deviation matrices, the time taken to compute the Sinkhorn projection is usually negligible.

For a doubly stochastic matrix, it is known that

$$0 \leq \log(\text{per}(nA)/n!) \leq n$$

the lower bound being attained at the matrix  $A = J$  whose components are all equal,  $J_{ij} = 1/n$ . The upper bound of  $\log(n^n/n!) \simeq n$  is attained at each of the permutation matrices, which are the extreme points in the set  $\text{DS}_n$  of doubly-stochastic matrices of order  $n$ . These bounds suggest that it may be easier to develop an approximation for the permanent of a doubly stochastic matrix than the permanent of an arbitrary positive matrix.

### 3 Moderate-deviation sequences

A sequence of square matrices  $X_1, X_2, \dots$  in which  $X_n$  is of order  $n$  is called *weakly bounded* if the absolute  $p$ th moment

$$\mu_p = \limsup_{n \rightarrow \infty} \frac{1}{n^2} \sum_{i,j=1}^n |X_n(i,j)|^p < \infty$$

is finite for all  $p \geq 1$ .

The set  $\text{DS}_n$  of doubly stochastic matrices of order  $n$  is convex. The extreme points are the  $n!$  permutation matrices, and the central point  $J$  is the equally-weighted average of the extremes. Given two matrices  $A, B \in \text{DS}_n$ , the row and column totals of the deviation matrix  $A - B$  are zero. It is helpful to define the  $L^2$ -norm and associated metric in  $\text{DS}_n$  by the standard Euclidean norm in the tangent space

$$d^2(A, B) = \sum_{i,j} (A_{ij} - B_{ij})^2, \quad \|A\|^2 = d^2(A, J) = \sum_{i,j} A_{ij}^2 - 1.$$

Thus, the central point has norm zero, and each extreme point has norm  $\sqrt{n-1}$ .

Each doubly stochastic matrix has one unit eigenvalue, and all others are less than or equal to one in absolute value. Consequently, it is helpful to define the operator norm in  $\text{DS}_n$  as the Hilbert-Schmidt norm of the deviation  $A - J$ , i.e. the largest absolute eigenvalue  $|\lambda_{\max}(A - J)|$ . The spectral gap of  $A$  is the difference  $\text{Gap}(A) = 1 - |\lambda_{\max}(A - J)|$ .

A sequence of doubly stochastic matrices  $\{A_n\}$  is said to be of *moderate deviation* if the re-scaled sequence  $X_n = n(A_n - J)$  is weakly bounded, and the spectral gap  $\text{Gap}(A_n) \geq C > 0$  is bounded below by a positive constant. If  $A, B$  are moderate-deviation sequences, the sequence  $AB$  of matrix products is also a moderate-deviation sequence, which implies that  $A'A$  is also of moderate deviation. Convex combinations are also of moderate deviation. Finally, for fixed  $\nu$ , the random sequence  $A_n \sim \text{DSD}_n(\nu)$  is of moderate deviation with probability one.

The main purpose of this note is to show that, for a moderate-deviation sequence

$$\log \text{per}(nA) = \log(n!) - \frac{1}{2} \log \det(I + J - A'A) + O(n^{-1})$$

with error decreasing as  $n \rightarrow \infty$ . One consequence is that  $\text{per}(nA)/n!$  is bounded as  $n \rightarrow \infty$ . Even though it does not affect the order of magnitude

Table 1: Numerical illustration of the determinantal approximation.

$n$	$\log \text{per}(K)$	$\rho = 1$		$\log \text{per}(K)$	$\rho = 2$	
		Approx	Error*		Approx	Error*
8	7.8905	7.8895	1.0178	5.7541	5.7454	8.7494
10	11.7601	11.7595	0.6188	9.0627	9.0577	5.0177
12	16.0163	16.0159	0.4319	12.7701	12.7667	3.3501
14	20.5955	20.5951	0.3293	16.8064	16.8039	2.4626
16	25.4522	25.4520	0.2636	21.1237	21.1218	1.9283
18	30.5526	30.5524	0.2199	25.6868	25.6852	1.5778
20	35.8700	35.8698	0.1880	30.4682	30.4668	1.3310
22	41.3831	41.3829	0.1636	35.4462	35.4451	1.1488
24	47.0743	47.0742	0.1534	40.6032	40.6022	1.0134

\*Error  $\times 10^3$

of the error, the modified approximation

$$\log \text{per}(nA) \simeq \log(n!) - \frac{1}{2} \log \det(I + t^2 J - t^2 A' A) \quad (1)$$

with  $t^2 = n/(n-1)$  is a worthwhile improvement for numerical work.

Table 1 illustrates the permanent approximation applied to a class of symmetric positive definite matrices  $K_{ij} = \exp(-\rho|x_i - x_j|)$  for points  $x_1, \dots, x_n$  equally spaced on the interval  $[0, 1]$ , so the components range from  $\exp(-\rho)$  to one. The first step uses the Sinkhorn algorithm to obtain a doubly stochastic matrix  $A$  from  $K$ , and the second step uses the determinantal approximation (1). All values shown are on the log scale. Because of the difficulty of computing the exact permanent, the range of  $n$ -values shown is rather limited. Nevertheless, it appears that the error decreases roughly as  $1/(300n)$  for  $\rho = 1$ , and  $1/(40n)$  for  $\rho = 2$ , in agreement with (1). A sequence of matrices of this type with bounded  $\rho$  is of moderate deviation; a sequence of matrices with  $\rho \propto n$  is not. It is unclear whether a sequence with  $\rho = \log(n)$  is moderate or not, but the approximation error does appear to decrease roughly at the rate  $1/n$  or  $\log(n)/n$ .

For a very special class of Kronecker-product matrices, Table 2 shows the relative error in the approximation for substantially larger values of  $n$ . These matrices have two blocks of size  $n/2$  each, with  $B_{ij} = 1$  within blocks and  $B_{ij} = \rho$  between blocks. The smaller the between-blocks value, the smaller the spectral gap, and the greater the approximation error. For the associated doubly stochastic matrix  $A = 2B/(n(1 + \rho))$ , Table 2 shows the exact value  $\log(\text{per}(nA)/n!)$  together with the error in the determinantal

Table 2: Determinantal approximation error for large  $n$ .

$n$	$\rho = 0.1$		$\rho = 0.05$		$\rho = 5/n$	
	Exact	Err $\times n$	Exact	Err $\times n$	Exact	Err
20	0.6755	1.6543	1.0187	2.8616	0.3177	0.0238
40	0.6274	1.3823	0.9616	3.4398	0.5366	0.0258
60	0.6140	1.2718	0.9274	3.1088	0.6900	0.0266
80	0.6082	1.2299	0.9114	2.8589	0.8076	0.0270
100	0.6049	1.2082	0.9029	2.7229	0.9029	0.0272
200	0.5987	1.1698	0.8882	2.5229	1.2155	0.0278
300	0.5967	1.1582	0.8839	2.4714	1.4065	0.0280
400	0.5957	1.1526	0.8818	2.4473	1.5444	0.0280

approximation  $-\log \det(I + J - A'A)/2$ . Once again, the empirical evidence suggests that the error for moderate-deviation matrices with constant  $\rho$  is  $O(n^{-1})$ , and for large-deviation matrices,  $O(1/(n\rho))$ .

## 4 Justification for the approximation

### 4.1 Permanent expansion

Let  $A$  be a matrix of order  $n$  such that  $nA_i^r = 1 + \epsilon_i^r$ , where the row and column totals of  $\epsilon$  are zero. If the components of  $\epsilon$  are real numbers greater than  $-1$ ,  $A$  is doubly stochastic and  $\epsilon/n = A - J$  is the deviation matrix. In the calculations that follow,  $\epsilon$  is complex-valued and weakly bounded, and  $\epsilon/n$  has spectral norm strictly less than one. In other words,  $A$  need not be doubly stochastic, or even real, but the associated sequence is assumed to be of moderate deviation.

The permanent expansion of  $nA$  as a polynomial of degree  $n$  has the form

$$\begin{aligned}
 \text{per}(1 + \epsilon) &= \sum_{\sigma} (1 + \epsilon_1^{\sigma(1)})(1 + \epsilon_2^{\sigma(2)}) \cdots (1 + \epsilon_n^{\sigma(n)}) \\
 &= n! \left( 1 + n \text{ave}(\epsilon_i^r) + \frac{n^{\downarrow 2}}{2!} \text{ave}^{\#}(\epsilon_i^r \epsilon_j^s) + \frac{n^{\downarrow 3}}{3!} \text{ave}^{\#}(\epsilon_i^r \epsilon_j^s \epsilon_k^t) + \cdots \right) \\
 &= n! \sum_{k=0}^n \frac{n^{\downarrow k}}{k!} \text{ave}^{\#}(\epsilon^{\otimes k}),
 \end{aligned}$$

where  $n^{\downarrow k} = n(n-1) \cdots (n-k+1)$  is the descending factorial, and  $\text{ave}^{\#}(\epsilon^{\otimes k})$  is the average of  $(n^{\downarrow k})^2$  products  $\epsilon_{i_1}^{r_1} \cdots \epsilon_{i_k}^{r_k}$  taken from distinct rows and

distinct columns. More explicitly, the term of degree  $k$  is

$$\frac{n^{\downarrow k}}{k!} \text{ave}^\#(\epsilon^{\otimes k}) = \frac{1}{k! n^{\downarrow k}} \sum^\# \epsilon_{i_1}^{r_1} \cdots \epsilon_{i_k}^{r_k}$$

with summation over distinct  $k$ -tuples  $(i_1, \dots, i_k)$  and  $(r_1, \dots, r_k)$ .

In the variational calculations that follow, each term such as the restricted sum  $\sum^\#(\epsilon^{\otimes k})$  or the unrestricted sum  $\sum(\epsilon^{\otimes k})$  is assigned a nominal order of magnitude as a function of  $n$ . For accounting purposes, the components of  $\epsilon$  are of order  $O(1)$ , and each product  $\epsilon_i^r \epsilon_j^s \epsilon_k^t$  is also regarded as being of order one, i.e. bounded, at least in a probabilistic sense, as  $n \rightarrow \infty$ . A sum such as  $\sum^\#(\epsilon^{\otimes k})$ , for fixed  $k$ , is of order  $O((n^{\downarrow k})^2) = O(n^{2k})$ , so the  $L^2$ -norm  $\sum(\epsilon_r^i)^2$  is of order  $O(n^2)$ . These sums could be of smaller order under suitable circumstances. For example, the difference  $(\sum - \sum^\#)\epsilon^{\otimes k}$  is a sum of  $n^{2k} - (n^{\downarrow k})^2$  terms, and is therefore of order  $O(n^{2k-2})$ . Thus, if the sum of the components of  $\epsilon$  is zero,  $\sum \epsilon^{\otimes k} = 0$  implies that the restricted sum  $\sum^\# \epsilon^{\otimes k}$  is of order  $O(n^{2k-2})$ . Likewise, if the components of  $\epsilon$  were independent random variables of zero mean,  $\sum \epsilon_{ir} = O(n)$ , and a similar conclusion holds for the restricted sum. The variational calculations that follow are not based on assumptions of statistical independence of components, but on the arithmetic implications of zero-sum constraints on rows and columns. For example,  $A$  may be symmetric.

The first goal is to show that the zero-sum restriction on the rows and columns of  $\epsilon$  implies that  $\sum^\#(\epsilon^{\otimes k})$  is of order  $O(n^k)$  rather than  $O(n^{2k})$ . In other words,  $\text{ave}^\#(\epsilon^{\otimes k})$  is of order  $O(n^{-k})$ . In fact, the even-degree terms in the permanent expansion are  $O(1)$ , while the odd terms are  $O(n^{-1})$ . These conclusions do not hold for all doubly stochastic matrices. For example, if  $A = (\rho I_n + J_n)/(1 + \rho)$  for some fixed  $\rho > 0$ , the  $L^2$ -norm  $\|A\| = \sqrt{n-1}\rho/(1+\rho)$  and other scalars of a similar type are not bounded.

## 4.2 Zero-sum restriction

In the term of degree three, the restricted sum is

$$\begin{aligned} \sum^\# \epsilon_i^r \epsilon_j^s \epsilon_k^t &= \sum^\# \epsilon_i^r \epsilon_j^s (\epsilon_i^r + \epsilon_i^s + \epsilon_j^r + \epsilon_j^s) \\ &= \sum^\# 2\epsilon_i^r \epsilon_j^s \epsilon_i^r + 2\epsilon_i^r \epsilon_j^s \epsilon_i^s \\ &= \sum^\# 2(\epsilon_i^r)^3 + 2(\epsilon_i^s)^3 \\ &= 4 \sum (\epsilon_i^r)^3. \end{aligned}$$

In the first line, the restricted sum over  $k \neq i, j$  is  $-\epsilon_i - \epsilon_j$ , while the restricted sum over  $t$  of  $-\epsilon_i^t$  is  $\epsilon_i^r + \epsilon_i^s$ . Proceeding in this way by restricted summation

over each non-repeated index, we arrive at the following expressions for the restricted tensorial sums of degree two, three and four:

$$\begin{aligned}
\sum^{\sharp} \epsilon^{\otimes 2} &= \sum^{\sharp} \epsilon_{i_1, i_2}^{r_1, r_2} = \sum \epsilon_{i_1, i_1}^{r_1, r_1} = \text{tr}(\epsilon' \epsilon) \\
\sum^{\sharp} \epsilon^{\otimes 3} &= 4 \sum^{\sharp} \epsilon_{i_1, i_1, i_1}^{r_1, r_1, r_1}, \\
\sum^{\sharp} \epsilon^{\otimes 4} &= 9 \sum^{\sharp} \epsilon_{i_1, i_1, i_1, i_1}^{r_1, r_1, r_1, r_1} - 9 \sum^{\sharp} \epsilon_{i_1, i_1, i_1, i_1}^{r_1, r_1, r_2, r_2} - 9 \sum^{\sharp} \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_1, r_1, r_1} \\
&\quad + 3 \sum^{\sharp} \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_1, r_2, r_2} + 6 \sum^{\sharp} \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_2, r_1, r_2} \\
&= 36 \sum \epsilon_{i_1, i_1, i_1, i_1}^{r_1, r_1, r_1, r_1} - 18 \sum \epsilon_{i_1, i_1, i_1, i_1}^{r_1, r_1, r_2, r_2} - 18 \sum \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_1, r_1, r_1} \\
&\quad + 3 \sum \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_1, r_2, r_2} + 6 \sum \epsilon_{i_1, i_1, i_2, i_2}^{r_1, r_2, r_1, r_2}.
\end{aligned}$$

Without the zero-sum constraint on the rows and columns,  $\sum^{\sharp} \epsilon^{\otimes k}$  is a sum over  $(n^{\downarrow k})^2$   $\epsilon$ -products, and thus of order  $O(n^{2k})$ . In the reduced form, each distinct value occurs in duplicate at least, so there are at most  $k/2$  distinct values for the row indices and  $k/2$  for the column indices. We observe that  $\sum^{\sharp} \epsilon^{\otimes 2} = \text{tr}(\epsilon' \epsilon)$  is  $O(n^2)$ , while  $\sum^{\sharp} \epsilon^{\otimes 3} = 4 \sum (\epsilon_i^r)^3$  is  $O(n^2)$  rather than  $O(n^3)$ . The five terms in  $\sum^{\sharp} \epsilon^{\otimes 4}$  are  $O(n^2)$ ,  $O(n^3)$ ,  $O(n^3)$ ,  $O(n^4)$  and  $O(n^4)$  respectively. The final pair can be expressed in matrix notation as  $3 \text{tr}^2(\epsilon' \epsilon) + 6 \text{tr}(\epsilon' \epsilon \epsilon' \epsilon)$ .

For any vector  $x = (x_1, \dots, x_k)$  in  $\mathfrak{R}^k$ , let  $\tau(x)$  be the associated partition of  $[k] = \{1, \dots, k\}$ , i.e.  $\tau(x)(r, s) = 1$  if  $x_r = x_s$  and zero otherwise. The restricted sum  $\sum^{\sharp} \epsilon^{\otimes k}$  can be expressed in reduced form either as a restricted sum or an unrestricted sum, the only difference arising in the coefficients as shown above for  $k = 4$ . Generally speaking, unrestricted sums are more convenient for computation. For general  $k$ , the restricted sums are as follows:

$$\begin{aligned}
\sum^{\sharp} \epsilon^{\otimes k} &= \sum_{\rho, \sigma} m^{\sharp}(\rho) m^{\sharp}(\sigma) \sum_{\substack{i: \tau(i) = \sigma \\ r: \tau(r) = \rho}} \epsilon_i^r \\
&= \sum_{\rho, \sigma} m(\rho) m(\sigma) \sum_{\substack{i: \tau(i) \geq \sigma \\ r: \tau(r) \geq \rho}} \epsilon_i^r,
\end{aligned}$$

where the outer sum extends over ordered pairs  $\rho, \sigma$  of partitions of the set  $[k]$ , and the inner sum over the row and column indices,  $i = (i_1, \dots, i_k)$  and  $r = (r_1, \dots, r_k)$ , which are held constant in each block.

The coefficients  $m, m^{\sharp}$  are multiplicative functions of the partition

$$m^{\sharp}(\rho) = \prod_{b \in \rho} (-1)^{\#b-1} (\#b - 1), \quad m(\rho) = \prod_{b \in \rho} (-1)^{\#b-1} (\#b - 1)!,$$



and  $m^\#(\rho) = m(\rho) = 0$  if  $\rho$  has a singleton block. Here, and elsewhere,  $\mathbf{1}_k$  denotes the maximal partition of  $[k]$  with one block,  $\#\rho$  is the number of blocks, and  $\#b$  is the number of elements of block  $b \in \rho$ . Although  $m$  and  $m^\#$  are related by Möbius inversion, these expressions are not to be confused with the Möbius function for the partition lattice:  $M(\rho, \mathbf{1}) = (-1)^{\#\rho-1}(\#\rho-1)!$ , which is a function of the number of blocks independently of their sizes.

For simplicity of notation in what follows, we write  $\epsilon_\sigma^\rho$  for the sum,

$$\epsilon_\sigma^\rho = \sum_{\substack{i:\tau(i)\geq\sigma \\ r:\tau(r)\geq\rho}} \epsilon_i^r,$$

in which  $i$  and  $r$  are constant within blocks, but the values for distinct blocks may be equal. Putting these expressions together with a scalar  $0 \leq t \leq 1$ , we find

$$\frac{\text{per}(1+t\epsilon)}{n!} = \sum_{k=0}^n \frac{t^k}{n \downarrow k k!} \sum_{\sigma, \rho \in \mathcal{P}_k} m(\rho)m(\sigma)\epsilon_\sigma^\rho \quad (2)$$

where  $\mathcal{P}_k$  is the set of partitions of  $[k]$ .

The polynomial (2) is the exponential generator  $\sum_{k=0}^n t^k \zeta_k / k!$  for the moment numbers

$$\zeta_k = \frac{1}{n \downarrow k} \sum_{\rho, \sigma \in \mathcal{P}_k} m(\rho)m(\sigma)\epsilon_\sigma^\rho$$

taking  $\zeta_k = 0$  for  $k > n$ . The coefficients in the log series are the cumulant numbers

$$\begin{aligned} \zeta'_k &= \sum_{\tau \in \mathcal{P}_k} (-1)^{\#\tau-1} (\#\tau-1)! \prod_{b \in \tau} \zeta_{\#b} \\ &= \sum_{\tau \in \mathcal{P}_k} (-1)^{\#\tau-1} (\#\tau-1)! \prod_{b \in \tau} \frac{1}{n \downarrow \#b} \sum_{\rho, \sigma \in \mathcal{P}_b} m(\rho)m(\sigma)\epsilon_\sigma^\rho \\ &= \sum_{\rho, \sigma \in \mathcal{P}_k} m(\rho)m(\sigma)\epsilon_\sigma^\rho \sum_{\tau \geq \rho \vee \sigma} (-1)^{\#\tau-1} (\#\tau-1)! \prod_{b \in \tau} \frac{1}{n \downarrow \#b} \\ &= \sum_{\rho, \sigma \in \mathcal{P}_k} m(\rho)m(\sigma)\epsilon_\sigma^\rho \times \Delta_n(\rho \vee \sigma) \end{aligned}$$

for  $k = 1, \dots$ . Note that  $\rho, \sigma$  in line 2 are the restrictions to the blocks of  $\tau$  of the partitions  $\rho, \sigma$  in line 3, so  $\tau \geq \rho, \sigma$ . Evidently,

$$\Delta_n(\rho) = \sum_{\sigma \geq \rho} (-1)^{\#\sigma-1} (\#\sigma-1)! \prod_{b \in \sigma} \frac{1}{n \downarrow \#b}$$

is the generalized cumulant associated with the reciprocal descending factorial series. For example,

$$\begin{aligned} n^{\downarrow 4} \Delta_n(12|34) &= 1 - \frac{n^{\downarrow 4}}{n^{\downarrow 2} n^{\downarrow 2}} = \frac{4n - 6}{n^{\downarrow 2}} \\ n^{\downarrow 5} \Delta_n(123|45) &= 1 - \frac{n^{\downarrow 5}}{n^{\downarrow 3} n^{\downarrow 2}} = \frac{6(n - 2)}{n^{\downarrow 2}} \\ n^{\downarrow 6} \Delta_n(12|34|56) &= 1 - \frac{3n^{\downarrow 6}}{n^{\downarrow 4} n^{\downarrow 2}} + \frac{2n^{\downarrow 6}}{(n^{\downarrow 2})^3} = \frac{8(n - 3)(7n - 10)}{(n^{\downarrow 2})^2}. \end{aligned}$$

Thus, the complete formal expansion for the log permanent is

$$\begin{aligned} \log(\text{per}(1 + t\epsilon)/n!) &= \sum_{k=1}^{\infty} \frac{t^k}{k!} \sum_{\rho, \sigma \in \mathcal{P}_k} m(\rho)m(\sigma)\epsilon_{\sigma}^{\rho} \Delta_n(\rho \vee \sigma) \\ &= \sum_{k=1}^{\infty} \frac{t^k}{k!} \sum_{\tau \in \mathcal{P}_k} \Delta_n(\tau) \sum_{\rho \vee \sigma = \tau} m(\rho)m(\sigma)\epsilon_{\sigma}^{\rho} \end{aligned}$$

in which  $\Delta_n(\mathbf{1}_k) = 1/n^{\downarrow k}$  for  $k \leq n$ , and zero otherwise.

### 4.3 Moderate-deviation asymptotic expansion

For a fixed pair of partitions  $\rho, \sigma \in \mathcal{P}_k$  the key scalar  $\epsilon_{\sigma}^{\rho}$  is the sum over indices  $\tau(i) \geq \sigma$  and  $\tau(r) \geq \rho$  of  $\epsilon_i^r$ . Under the rules for determining the order of magnitude in  $n$ ,  $\epsilon_{\sigma}^{\rho}$  is of order  $O(n^{\#\sigma + \#\rho})$ , and

$$\Delta_n(\rho \vee \sigma) \epsilon_{\sigma}^{\rho} = O(n^{\#\sigma + \#\rho - k - \#(\rho \vee \sigma) + 1}).$$

Since neither partition has singleton blocks,  $2\#\sigma \leq k$  and  $2\#\rho \leq k$ . Thus the order is  $O(1)$  only if  $k$  is even, each partition is binary with all blocks of size two, and the least upper bound is the full set  $\mathbf{1}_k$ . Otherwise, if the least upper bound is less than  $\mathbf{1}_k$ , or if any block of either partition is of size three or more, the term is  $O(n^{-1})$  or smaller.

The leading term of maximal order in the expansion of the log permanent is

$$\begin{aligned} h_n^{(0)}(t) &= \sum_{k=1}^{n/2} \frac{t^{2k}}{(2k)! n^{\downarrow 2k}} \sum_{\substack{\rho, \sigma \sim 2^k \\ \rho \vee \sigma = \mathbf{1}_{2k}}} \epsilon_{\sigma}^{\rho} \\ &= \sum_{k=1}^{n/2} \frac{t^{2k}}{2k} \text{tr}((\epsilon' \epsilon)^k) / n^{\downarrow 2k} \end{aligned}$$

$$\begin{aligned}
&= \sum_{k=1}^{\infty} \frac{t^{2k}}{2k} \operatorname{tr}((\epsilon' \epsilon / n^{\downarrow 2})^k) + O(n^{-1}) \\
&= -\frac{1}{2} \log \det(I_n - t^2 \epsilon' \epsilon / n^{\downarrow 2}) + O(n^{-1}) \\
h_n^{(0)}(1) &= -\frac{1}{2} \log \det(I_n + J - A' A) + O(n^{-1}).
\end{aligned}$$

The symbol  $\rho, \sigma \sim 2^k$  denotes two partitions of  $[2k]$  having  $k$  blocks of size two, and there are  $(2k-1)!$  pairs whose least upper bound is the full set  $[2k]$ . The series is convergent for  $|t| < 1/|\lambda_{\max}(A - J)|$ , so  $h_n^{(0)}(1)$  is finite by the spectral gap assumption.

In the expansion of the first-order correction  $h^{(1)}(t)$ , the terms may be grouped by degree in  $t$ . The following are the terms of degree six or less, expressed so far as possible using matrix operations in which  $\epsilon_r$  and  $(\epsilon' \epsilon)_r$  are the component-wise  $r$ th powers of  $\epsilon$  and  $(\epsilon' \epsilon)$  respectively.

$$\begin{aligned}
\text{degree 3: } & \frac{2}{3n^3} \sum \epsilon_3 = \frac{2}{3n^3} \sum (\epsilon_i^r)^3 \\
\text{degree 4: } & -\frac{3}{4n^4} \sum (\epsilon'_2 \epsilon_2 + \epsilon_2 \epsilon'_2) \\
\text{degree 5: } & \frac{2}{n^5} \operatorname{tr}(\epsilon_2 \epsilon' \epsilon \epsilon') + \frac{1}{n^5} \sum \epsilon_2 \epsilon' \epsilon_2 \\
\text{degree 6: } & \frac{1}{3n^6} \sum ((\epsilon' \epsilon)_3 + (\epsilon \epsilon')_3) + \frac{1}{2n^6} \sum (\epsilon'_2 \epsilon \epsilon' \epsilon_2 + \epsilon_2 \epsilon' \epsilon \epsilon'_2) \\
& -\frac{3}{2n^6} \sum (\epsilon_2 (\epsilon' \epsilon)_2 + \epsilon'_2 (\epsilon \epsilon')_2)
\end{aligned}$$

## 5 Random matrices

In order to test the adequacy of the determinantal approximation, 5000 random matrices  $A$  of order 10–25 were generated, and the exact values  $y(A) = \log(\operatorname{per}(nA)/n!)$  were computed using the Ryser algorithm. The matrices of even order were generated from the distribution  $A_i \sim \operatorname{DSD}_n(\nu_i)$ . In order to ensure an adequate range of permanent values, the parameters  $\nu_i$  were generated randomly and independently from the Gamma distribution with mean one on four degrees of freedom, i.e.  $\chi_8^2/8$ . On the matrices of odd order was superimposed a multiplicative random block pattern  $(\eta + \eta' B(r, s))$  in which  $\eta, \eta'$  are uniform  $(0, 1)$  random scalars, and  $B$  is a random partition of  $[n]$ , the Chinese restaurant process with parameter 1. This block pattern was superimposed prior to Sinkhorn projection in order to vary the eigenvalue pattern, which depends on the coefficients  $\eta$  and on the block sizes. Matrices of this type do not satisfy the moderate-deviation condition.

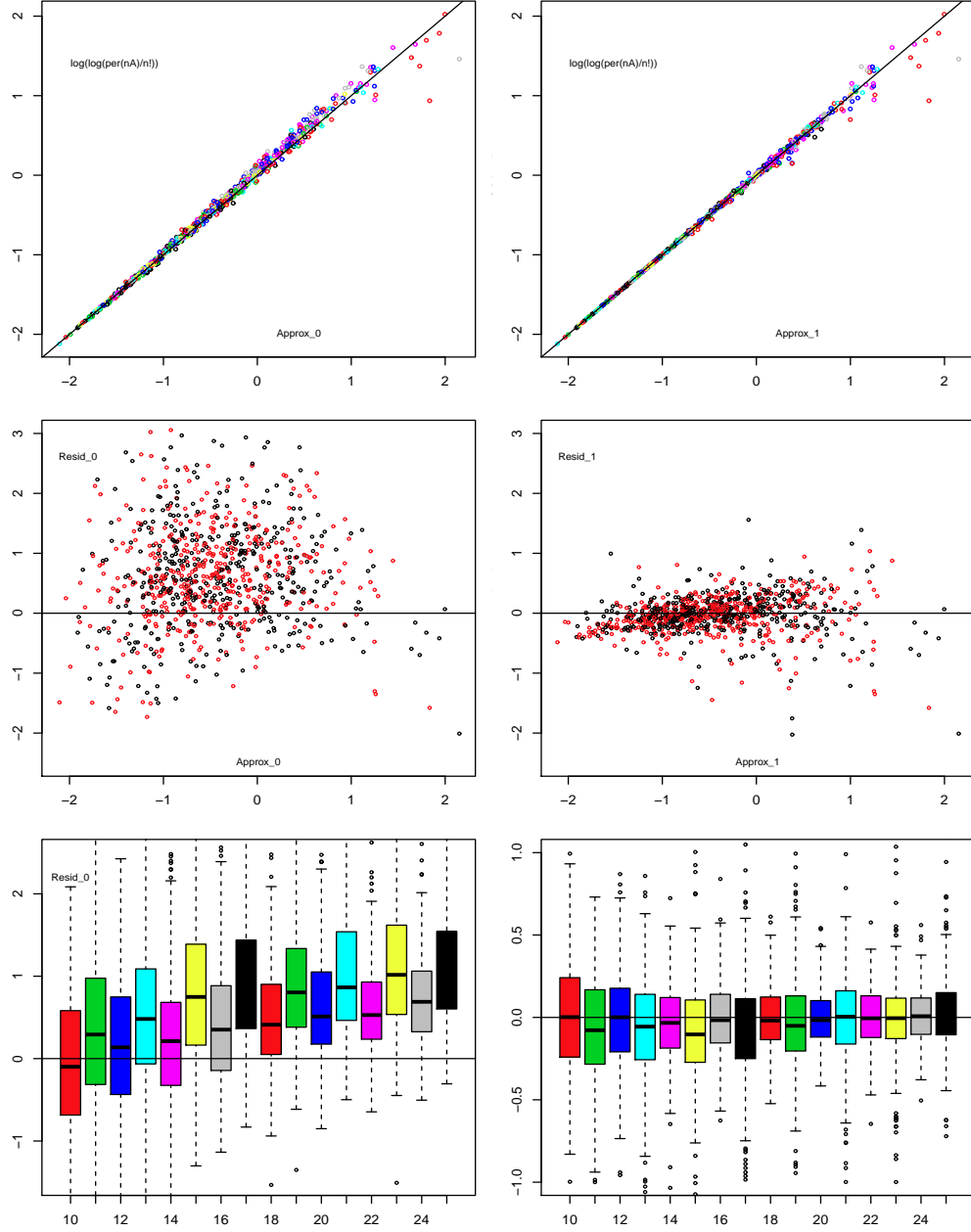


Figure 1. Top row: exact permanent of 500 random matrices plotted against two approximations, both on the log-log scale. Middle row: standardized log-log residuals plotted against the two approximations. Third row: residuals plotted against  $n$ , at twice the scale in the right panel.

On the log scale, the range of observed permanental values was  $0.11 < y(A) < 7.56$ , or  $0.042 < y/\log(n) < 2.57$ , the largest value occurring for a matrix of order 19. The target range  $y(A) \leq 2$  was exceeded by 2.6% of the matrices generated, and the moderate-deviation threshold  $y > \log(n)$  was exceeded by 43 matrices comprising roughly 0.9% of the simulations. Most of these exceedances occurred for matrices of odd order having a pronounced block structure. The range of  $L^2$ -values was  $0.22 \leq \|A\|^2 \leq 7.32$ , with only 2% in excess of 3.0. Although there exist matrices  $A \in \text{DS}_n$  such that the ratio  $y(A)/\|A\|^2$  is arbitrarily large or arbitrarily small, the range of simulated values was only  $(0.53, 1.15)$ . This is a reminder that the behaviour of the permanent in a bounded  $L^2$ -ball is not a good indicator of its behaviour near the corners.

Two approximations were computed, the first-order determinantal approximation  $x_0(A)$  as in (1), and a second-order modification  $x_1(A)$  using the additional terms in section 4.3. The scatterplots of  $\log \log y(A)$  against  $\log \log x_0(A)$  and  $\log \log x_1(A)$  are shown in the top two panels of Fig. 1. To reduce clutter in the top two rows, only a 10% sample is shown.

Since the relative error of the approximation  $x(A)$  increases with its magnitude, the standardized residuals are most naturally defined on the log-log scale

$$\text{Resid}(y, x) = \frac{n \log(y(A)/x(A))}{x(A)}.$$

In the middle panels of Fig. 1, the log-log residuals are plotted against  $x$ , using the same scale for both plots. In the lower panels, the log-log residuals are plotted against  $n$ . The first plot makes it clear that the determinantal approximation  $x_0(A)$  tends to be an under-estimate: the rate of occurrence of the inequality  $y(A) > x_0(A)$  increases from 44% for  $n = 10$  to over 90% for  $n \geq 20$ . In the lower left panel, the alternating pattern of residuals for odd and even  $n$  is a consequence of the block pattern embedded in the matrices of odd order. For both types of matrices, the plots suggest that the residual distribution is asymptotically constant as  $n \rightarrow \infty$ , and that the correction term is appreciable for moderate  $n$ . The root-mean-squared residual is approximately  $1.2/n$  for the first approximation and  $0.3/n$  for the second, but the distributions are more like Cauchy than Gaussian. As it happens, the inequalities

$$\log y(A) < \log x_0(A) - 1.4 x_0(A)/n, \quad \log y(A) > \log x_0(A) + 3.0 x_0(A)/n$$

occur in the sample with rates 1.0% each. The inequality

$$|\log y(A) - \log x_1(A)| > 0.5 x_1(A)/n$$

occurs at a fairly constant rate of around 1% for non-block-structured matrices. For block-structured matrices, the rate is about five times as high, but slowly decreasing in  $n$  over the range observed. Of course, these rates must depend on the distribution by which the matrices are generated, but it is anticipated that the rate will have a positive limit for matrices in the moderate-deviation region  $x_0(A) < \log(n)$ . In other words, we should expect  $\log(\text{per}(nA)/n!)$  to lie in the interval  $x_1(A) \exp(\pm x_1(A) \log(n)/n)$  with probability tending to one for large  $n$  if  $x_1(A)$  is bounded.

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